

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034524.D
 Acq On : 10 Oct 2024 16:55
 Operator : RC/JU
 Sample : P4263-06
 Misc :
 ALS Vial : 100 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE114D2-20240930

Quant Time: Oct 10 17:51:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 26 14:44:02 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	4252	0.400	ng	-0.02
7) Naphthalene-d8	10.127	136	12054	0.400	ng	-0.02
13) Acenaphthene-d10	14.019	164	6850	0.400	ng	-0.03
19) Phenanthrene-d10	16.784	188	13020	0.400	ng	#-0.02
29) Chrysene-d12	21.008	240	7655	0.400	ng	0.00
35) Perylene-d12	23.116	264	10465	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	1613	0.134	ng	-0.01
5) Phenol-d6	6.600	99	589	0.042	ng	0.00
8) Nitrobenzene-d5	8.525	82	2428	0.263	ng	-0.01
11) 2-Methylnaphthalene-d10	11.731	152	6299	0.354	ng	-0.02
14) 2,4,6-Tribromophenol	15.542	330	923	0.297	ng	0.00
15) 2-Fluorobiphenyl	12.640	172	8911	0.320	ng	-0.02
27) Fluoranthene-d10	18.829	212	13718	0.418	ng	-0.02
31) Terphenyl-d14	19.447	244	11027	0.721	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	3.025	88	19993	3.761	ng	Qvalue 99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
Data File : BN034524.D
Acq On : 10 Oct 2024 16:55
Operator : RC/JU
Sample : P4263-06
Misc :
ALS Vial : 100 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
RE114D2-20240930

Quant Time: Oct 10 17:51:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 26 14:44:02 2024
Response via : Initial Calibration

